

Data Path : Z:\HPCHEM1\BNA E\DATA\BE062416\
 Data File : BE091944.D
 Acq On : 24 Jun 2016 13:34
 Operator : UM/SJ
 Sample : H3676-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 NCHC-008

Quant Time: Jun 24 17:26:43 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE061416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 15 12:00:38 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.20	152	30449	5.00	ng	0.00
8) Naphthalene-d8	11.02	136	142574	5.00	ng	0.02
15) Acenaphthene-d10	14.86	164	106168	5.00	ng	0.00
24) Phenanthrene-d10	17.61	188	202418	5.00	ng	0.00
31) Chrysene-d12	21.76	240	289621	5.00	ng	0.00
40) Perylene-d12	24.17	264	212218	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.72	112	41869	7.03	ng	0.00
5) Phenol-d6	7.37	99	63916	6.56	ng	0.02
9) Nitrobenzene-d5	9.36	82	35460	3.60	ng	0.00
16) 2,4,6-Tribromophenol	16.35	330	17429	6.32	ng	0.00
17) 2-Fluorobiphenyl	13.48	172	91046	3.38	ng	0.00
34) Terphenyl-d14	20.22	244	135469	3.82	ng	0.00
Target Compounds						
27) Pentachlorophenol	17.26	266	92	0.11	ng	# 82
30) Fluoranthene	19.65	202	35462	0.16	ng	# 81
33) Pyrene	20.02	202	34584	0.18	ng	# 68
35) Benzo(a)anthracene	21.73	228	5107m	0.11	ng	
37) Chrysene	21.78	228	6205m	0.09	ng	
39) Indeno(1,2,3-cd)pyrene	26.68	276	16484	0.07	ng	# 87
41) Benzo(b)fluoranthene	23.43	252	7622m	0.14	ng	
42) Benzo(k)fluoranthene	23.46	252	3015m	0.06	ng	
43) Benzo(a)pyrene	24.06	252	4744	0.10	ng	# 25
45) Benzo(g,h,i)perylene	27.47	276	16728	0.09	ng	# 60

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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